

Figure 1: Association [3H]-4MG

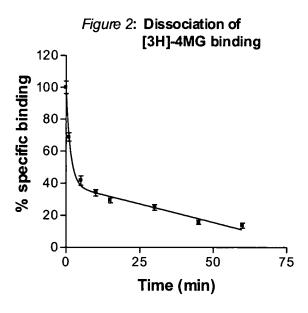


Figure 3: DRUG INHIBITION OF [3H]-4MG BINDING

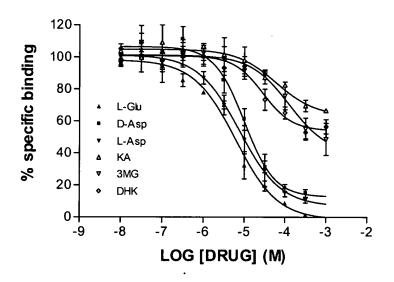
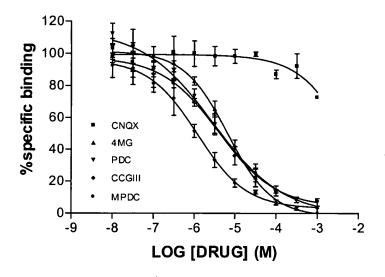


Figure 4: DRUG INHIBITION of [3H]-4MG BINDING



% specific binding

120 - 100 -

NMDA

4MG

MK801(1mM) GLY (1mM)

GABA (1mM)

Figure 5: COM PARISON of MISCELLANEOUS DRUG INHIBITION of [3H]-D-ASPARTATE and [3H]-4MG

AMPA

АТРА FW D-Asp

$$NH_2$$
 HO_2C
 $CONHSO_2CH_3$
 CH_3O_2SHNOC
 CO_2H
 CO_2H
 $CO_2CONHSO_2CF_3$

Figure 6A

Figure 6B

Figure 6C

HO PhHNCO2 NHCONHPh
$$NH_2$$
 NH_2 N

$$H_3CO_2C$$
 NH_2
 HO_2C
 CO_2H
 NH_2
 N

NHCONHPh
$$HO_2C$$

$$OH$$

$$HO_2C$$

$$CO_2H$$

$$OH$$

$$HO_2C$$

$$CO_2H$$

$$HO_2C$$

$$CO_2H$$

Figure 6D

Figure 6E

Figure 6F

со⁵н

Figure 6G

CO₂H

HO₂C



$$CO_2H$$
 NH_2
 NH_2

MeO₂C

$$\begin{array}{c}
 & \text{NH}_2 \\
 & \text{HO}_2C
\end{array}$$
 $\begin{array}{c}
 & \text{NH}_2 \\
 & \text{HO}_2C
\end{array}$
 $\begin{array}{c}
 & \text{NH}_2 \\
 & \text{HO}_2C
\end{array}$

Figure 6H

$$RO_2C$$
 $R = Me, Et, {}^{t}Bu$
 R^{1}
 R^{2}
 R^{2}
 R^{3}
 R^{4}
 R^{5}
 R^{5}

R¹ = CH₃, and halogen

R², R³ are independently

H, C1-C6-alkyl, C3-C4-alkenyl, C3-C5-cycloalkyl, C1-C6-alkyl-CO-,

C1-C6-alkyl-OCO-, C1-C6-alkyl-NHCO-, HCO-, or C3-C6-alkynyl

 R^2 , R^3 taken together can be $-CH_2(CH_2)_pCH_2$ -

$$R$$
 CO_2H
 HO_2C
 $R = H, Me, Et, Cl$
 $R = H, Me, Et, nPr$
 $R = H, Et, nPr$